LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

Signal : TIC: BP021117.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.187	30	34	41	rVB	24223	35149	0.60%	0.057%
2	3.705	118	122	131	rVB	36297	55912	0.95%	0.090%
3	3.911	153	157	162	rVB	25759	33722	0.57%	0.054%
4	4.322	224	227	236	rVB	9691	19654	0.33%	0.032%
5	4.811	306	310	318	rVB	44783	69462	1.18%	0.112%
6	5.681	453	458	468	rBV2	15591	31109	0.53%	0.050%
7	6.764	635	642	650	rBV2	31361	54613	0.93%	0.088%
8	6.881	657	662	678	rVV	326490	635625	10.82%	1.027%
9	7.011	678	684	695	rVV	299681	492227	8.38%	0.795%
10	7.216	713	719	730	rBV	386390	667422	11.36%	1.079%
11	7.299	730	733	740	rVB3	10745	22716	0.39%	0.037%
12	7.664	787	795	805	rBV	656980	1016616	17.30%	1.643%
13	8.399	912	920	930	rBV	417336	746508	12.70%	1.206%
14	8.493	932	936	946	rVB2	27488	61375	1.04%	0.099%
15	8.822	985	992	1008	rBV	371210	664994	11.32%	1.075%
16	9.163	1044	1050	1057	rVB2	21247	35136	0.60%	0.057%
17	9.375	1075	1086	1093	rBV3	37470	81504	1.39%	0.132%
18	9.534	1105	1113	1125	rBV	289206	540838	9.20%	0.874%
19	10.093	1199	1208	1226	rBV	400858	866768	14.75%	1.401%
20	10.434	1259	1266	1272	rBV	880045	1462235	24.88%	2.363%
21	10.481	1272	1274	1282	rVB	35532	53421	0.91%	0.086%
22	10.599	1286	1294	1307	rBV	166862	387963	6.60%	0.627%
23	10.746	1314	1319	1327	rVB9	11453	23494	0.40%	0.038%
24	10.981	1355	1359	1369	rBV6	14443	30253	0.51%	0.049%
25	11.187	1387	1394	1404	rBV2	125573	296178	5.04%	0.479%
26	11.652	1468	1473	1479	rBV2	13714	21934	0.37%	0.035%
27	12.110	1545	1551	1556	rBV	28043	52727	0.90%	0.085%
28	12.322	1582	1587	1594	rVB2	27415	47953	0.82%	0.077%
29	12.475	1609	1613	1619	rBV4	10620	22329	0.38%	0.036%
30	12.746	1651	1659	1667	rBV4	8384	20445	0.35%	0.033%
31	13.099	1713	1719	1722	rBV5	14930	27961	0.48%	0.045%
32	13.134	1722	1725	1729	rVV2	12348	18521	0.32%	0.030%
33	13.187	1729	1734	1738	rVB	13441	24684	0.42%	0.040%
34	13.616	1801	1807	1812	rBV5	21071	43485	0.74%	0.070%
35	13.675	1812	1817	1818	rBV3	15552	20209	0.34%	0.033%
36	13.710	1818	1823	1830	rVV	816169	1196959	20.37%	1.934%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

37	14.004	1861	1873	1883	rBV	1022396	1725038	29.35%	2.788%
38	14.128	1889	1894	1900	rVB5	23134	46869	0.80%	0.076%
39	14.187	1900	1904	1911	rVB5	10296	17784	0.30%	0.029%
40	14.310	1917	1925	1931	rBV	1243106	2053574	34.94%	3.319%
41	14.375	1931	1936	1943	rVB	119113	201064	3.42%	0.325%
42	14.504	1952	1958	1960	rBV2	433830	657352	11.19%	1.062%
43	14.616	1971	1977	1990	rBV4	109533	337531	5.74%	0.545%
44	14.722	1990	1995	1999	rVV	71632	144622	2.46%	0.234%
45	14.769	1999	2003	2010	rVB5	45457	86237	1.47%	0.139%
46	14.987	2035	2040	2045	rVV8	7261	17736	0.30%	0.029%
47	15.104	2055	2060	2064	rBV6	16419	29827	0.51%	0.048%
48	15.151	2064	2068	2074	rVV2	23411	41915	0.71%	0.068%
49	15.322	2091	2097	2103	rBV	1461796	2003521	34.09%	3.238%
50	15.381	2103	2107	2112	rVB2	130313	192358	3.27%	0.311%
51	15.493	2121	2126	2132	rBV	159201	244752	4.16%	0.396%
52	15.681	2153	2158	2164	rVB2	23789	43473	0.74%	0.070%
53	15.834	2181	2184	2190	rBV	26735	43066	0.73%	0.070%
54	15.892	2190	2194	2202	rVB6	32678	65107	1.11%	0.105%
55	16.081	2217	2226	2234	rBV8	22713	74065	1.26%	0.120%
56	16.287	2257	2261	2266	rBV3	14023	21794	0.37%	0.035%
57	16.410	2276	2282	2286	rBV4	35943	60814	1.03%	0.098%
58	16.463	2287	2291	2296	rVB5	20015	30493	0.52%	0.049%
59	16.528	2298	2302	2305	rBV5	17693	26338	0.45%	0.043%
60	16.645	2318	2322	2325	rBV5	17556	25080	0.43%	0.041%
61	16.704	2330	2332	2339	rVB	53920	76820	1.31%	0.124%
62	16.781	2341	2345	2352	rVB	57330	86730	1.48%	0.140%
63	16.869	2352	2360	2365	rBV5	46257	119596	2.04%	0.193%
64	16.939	2368	2372	2377	rVB	83436	124916	2.13%	0.202%
65	17.128	2397	2404	2407	rBV	1652360	2523821	42.94%	4.079%
66	17.169	2407	2411	2416	rVV	1576171	2440912	41.53%	3.945%
67	17.228	2416	2421	2426	rVV2	1337731	2189249	37.25%	3.538%
68	17.263	2426	2427	2432	rVB	309634	347571	5.91%	0.562%
69	17.557	2468	2477	2486	rBV2	145423	370181	6.30%	0.598%
70	17.657	2492	2494	2500	rBV	24684	34126	0.58%	0.055%
71	17.804	2516	2519	2523	rBV	56807	70400	1.20%	0.114%
72	18.028	2553	2557	2558	rBV	169919	187892	3.20%	0.304%
73	18.057	2558	2562	2564	rVV	423141	611012	10.40%	0.987%
74	18.086	2564	2567	2571	rVB	263260	372040	6.33%	0.601%
75	18.163	2578	2580	2585	rVB	82201	94683	1.61%	0.153%
76	18.228	2586	2591	2595	rVV3	355744	607449	10.34%	0.982%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Title : SVOA CALIBRATION

77	18.263	2595	2597	2606	rVB	137157	207777	3.54%	0.336%
78	18.410	2619	2622	2625	rBV4	36835	54080	0.92%	0.087%
79	18.545	2641	2645	2650	rVV	151925	200343	3.41%	0.324%
80	18.604	2651	2655	2664	rVB	202826	348939	5.94%	0.564%
81	19.034	2725	2728	2733	rVB4	129386	192600	3.28%	0.311%
82	19.145	2745	2747	2753	rVB2	99631	130257	2.22%	0.210%
83	19.263	2761	2767	2773	rBV	3653720	4764766	81.08%	7.700%
84	19.386	2785	2788	2796	rVB3	209889	362708	6.17%	0.586%
85	19.610	2820	2826	2828	rBV2	2163021	3184699	54.19%	5.147%
86	19.639	2828	2831	2839	rVB	2488840	3361025	57.19%	5.431%
87	20.151	2914	2918	2923	rVB2	424948	705378	12.00%	1.140%
88	20.257	2933	2936	2940	rVB	277606	296554	5.05%	0.479%
89	21.128	3081	3084	3088	rBV2	247083	368888	6.28%	0.596%
90	21.175	3088	3092	3094	rVV2	420484	592126	10.08%	0.957%
91	21.210	3094	3098	3105	rVB4	761205	1349818	22.97%	2.181%
92	21.445	3135	3138	3144	rVB	423579	605377	10.30%	0.978%
93	21.557	3149	3157	3162	rVV2	2273445	5876890	100.00%	9.497%
94	21.604	3162	3165	3179	rVB	1419155	2535490	43.14%	4.097%
95	21.763	3188	3192	3197	rVB2	331019	519832	8.85%	0.840%
96	21.839	3202	3205	3211	rVB2	265460	390797	6.65%	0.632%
97	23.827	3538	3543	3550	rBV	814578	2016096	34.31%	3.258%
98	24.651	3678	3683	3689	rVV	595785	1279182	21.77%	2.067%
99	24.716	3690	3694	3704	rVB	570612	1536903	26.15%	2.484%
100	24.874	3716	3721	3734	rVB	1006211	2669955	45.43%	4.315%

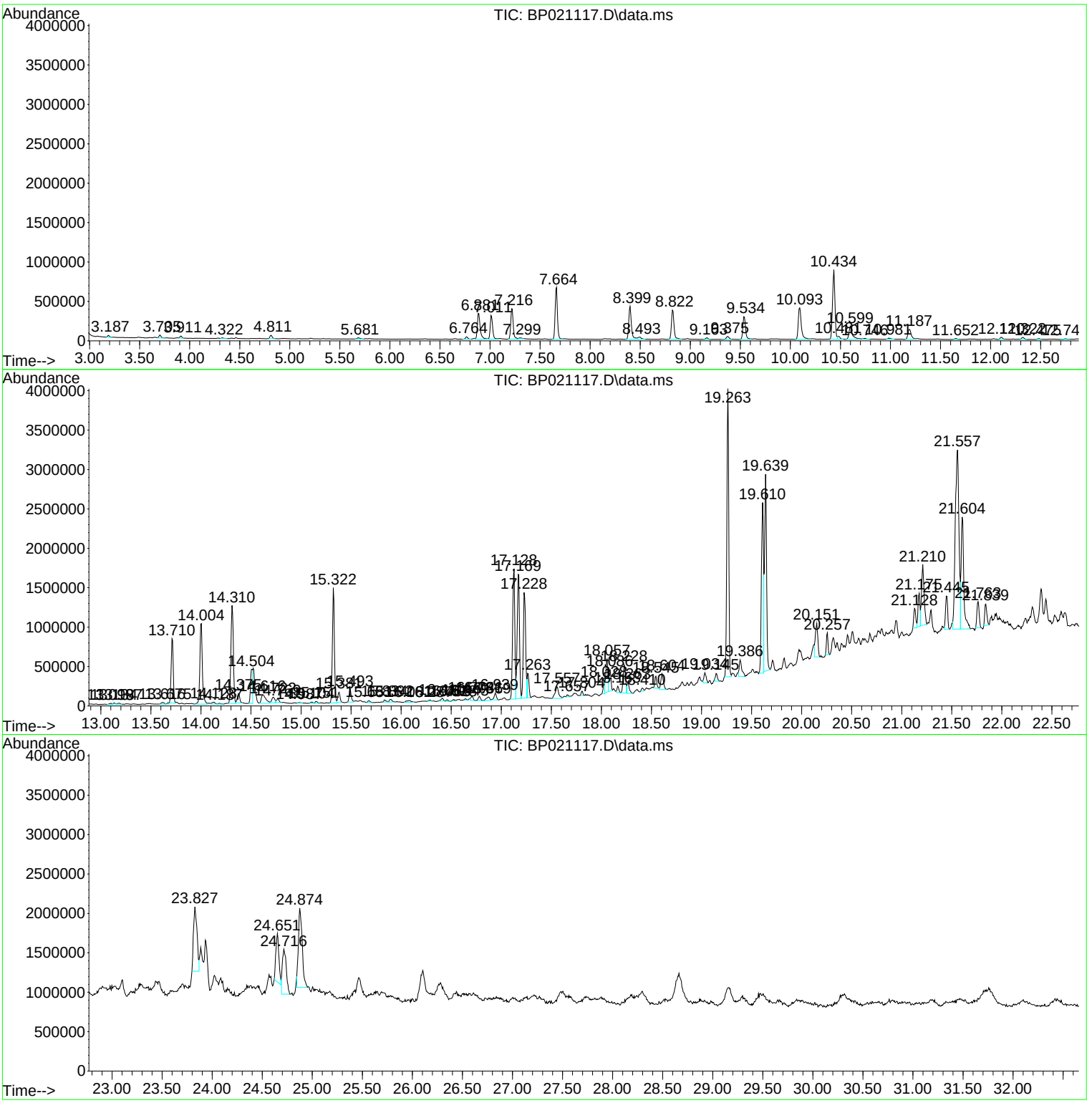
Sum of corrected areas: 61880389

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

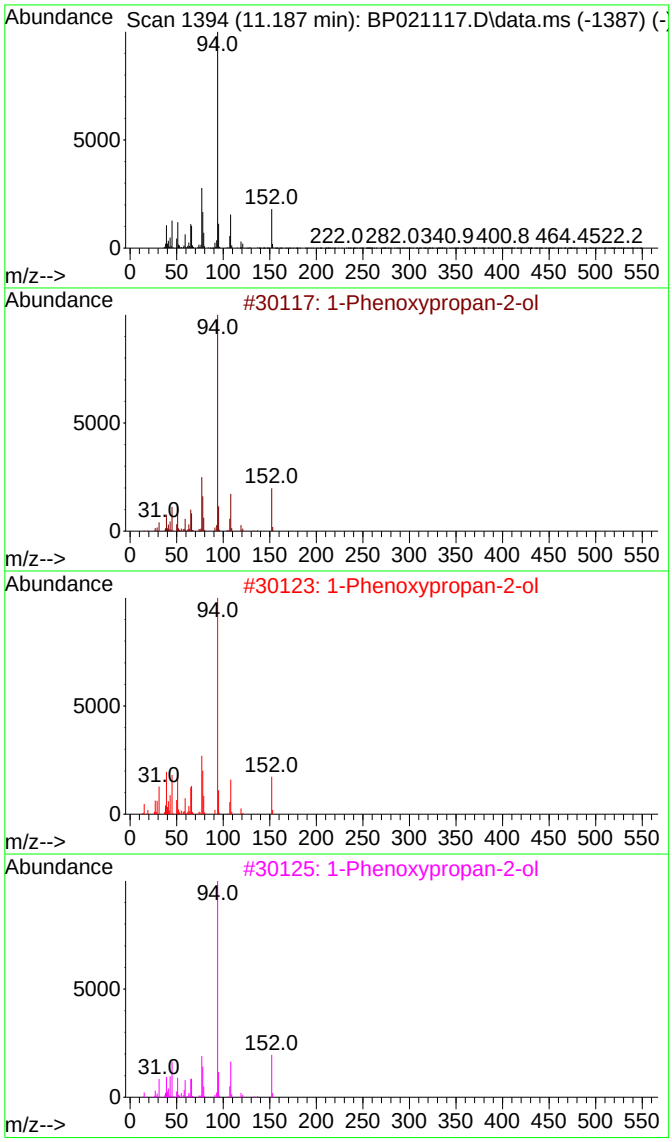
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

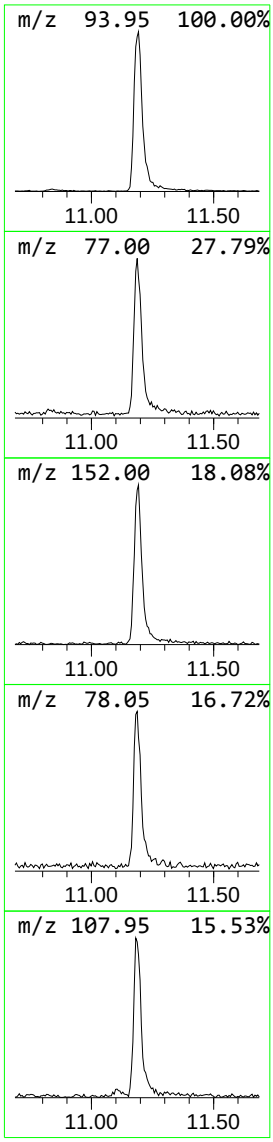
****

Peak Number 1 1-Phenoxypropan-2-ol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.187	4.05 ng/ul	296178	Naphthalene-d8	10.434

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	96
2	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	94
3	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
4	1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
5	1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

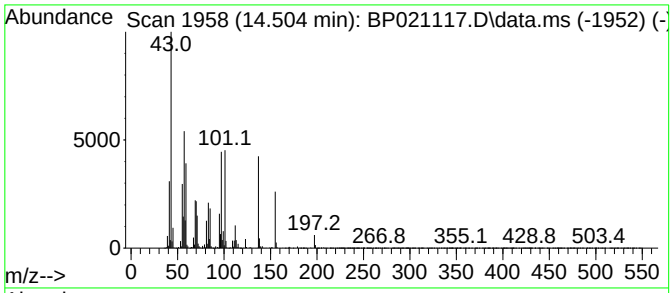
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

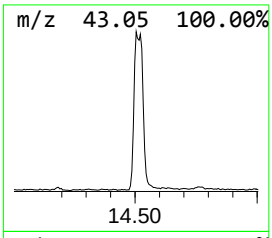
****

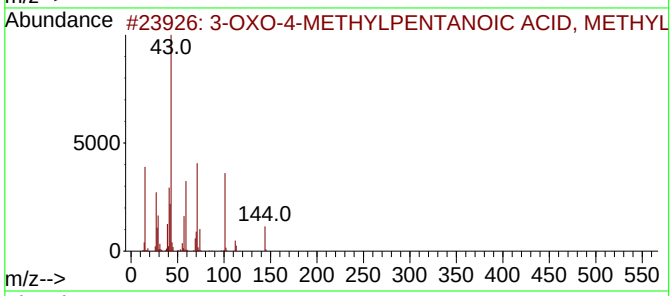
Peak Number 2 unknown-01 Concentration Rank 1

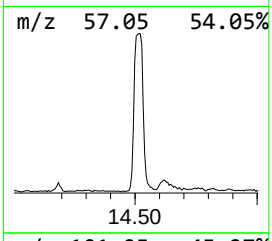
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.504	6.40 ng/ul	657352	Acenaphthene-d10	14.310

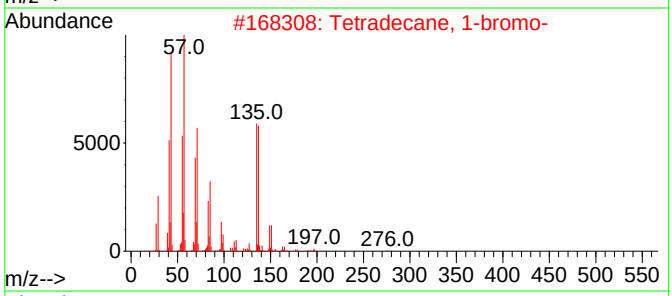
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-OXO-4-METHYLPENTANOIC ACID, ME...	144	C7H12O3	042558-54-3	12
2	Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	10
3	Carbonic acid, prop-1-en-2-yl un...	256	C15H28O3	1000382-90-6	10
4	Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	10
5	2-Pentanol, 2-methyl-, acetate	144	C8H16O2	034859-98-8	10

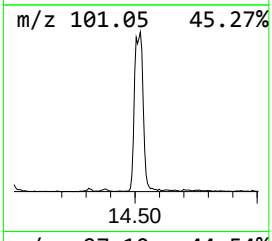


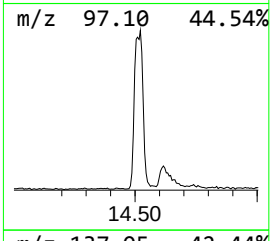


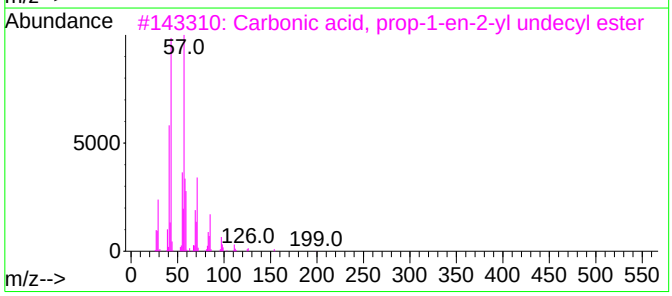


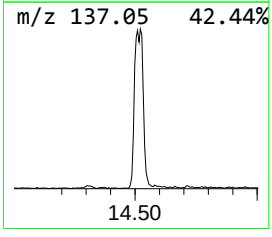












Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

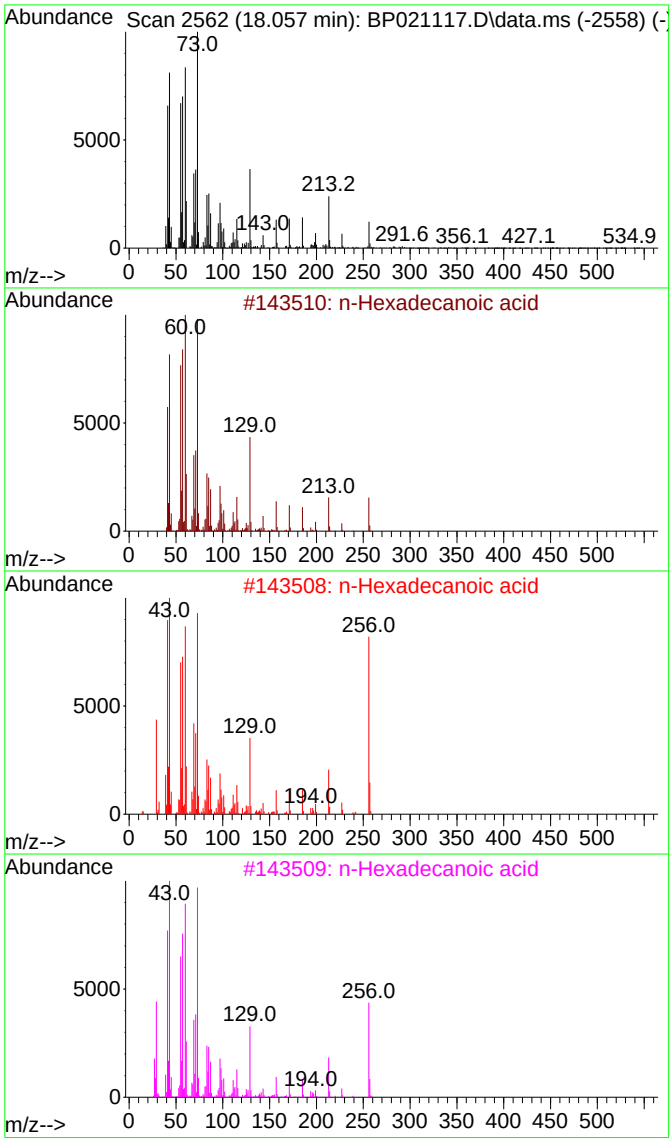
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

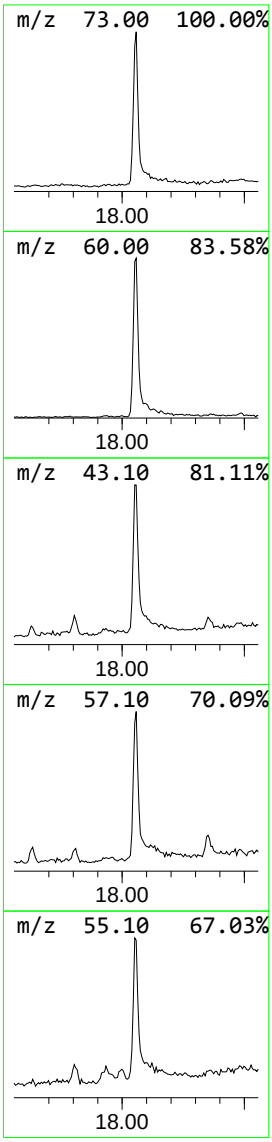
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.057	4.84 ng/ul	611012	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

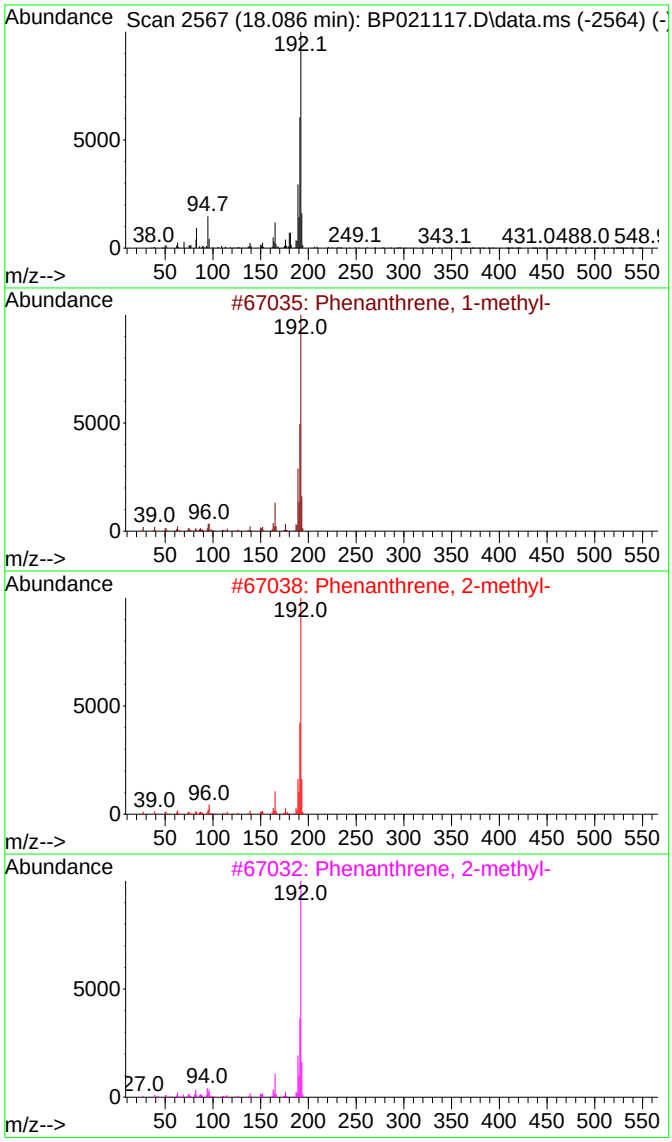
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

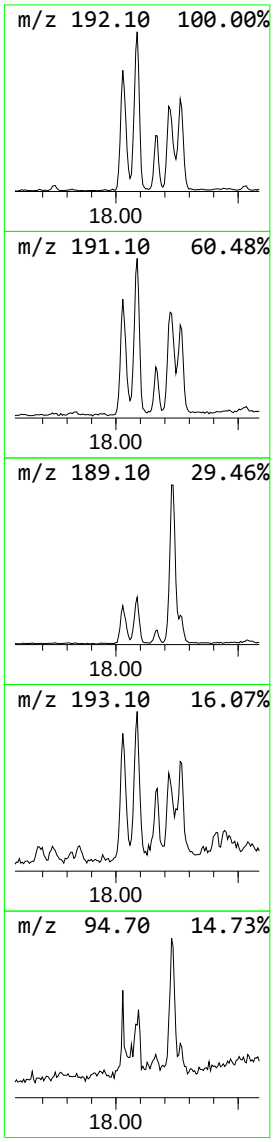
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Phenanthrene, 1-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.087	2.95 ng/ul	372040	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	97
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	96
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	95
4	1H-Cyclopropa[l]phenanthrene,1a,...	192	C15H12	000949-41-7	94
5	Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	93





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Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

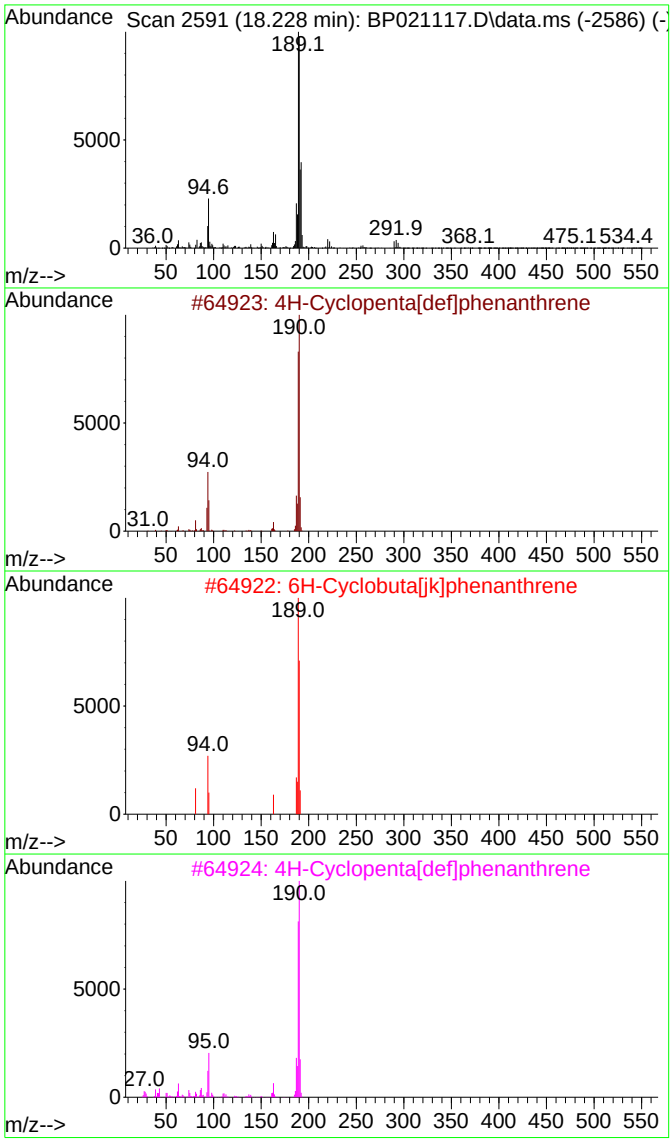
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

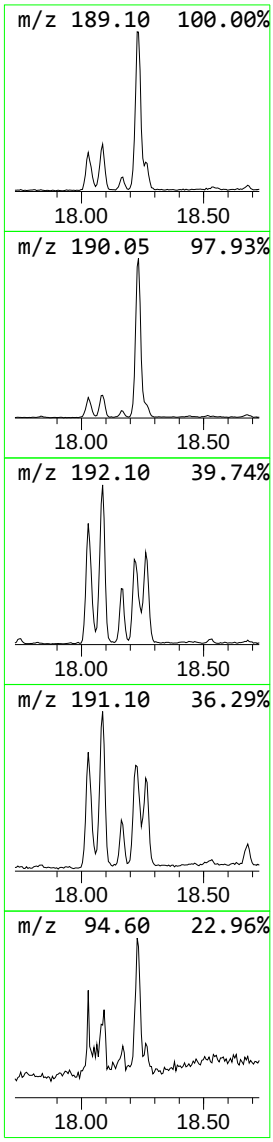
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 4H-Cyclopenta[def]phenanthrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.228	4.81 ng/ul	607449	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	93
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	59
3	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	58
4	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	52
5	1H-Pyrazole-4-carboxaldehyde, 1-...	190	C10H7FN2O	1000338-05-6	47





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

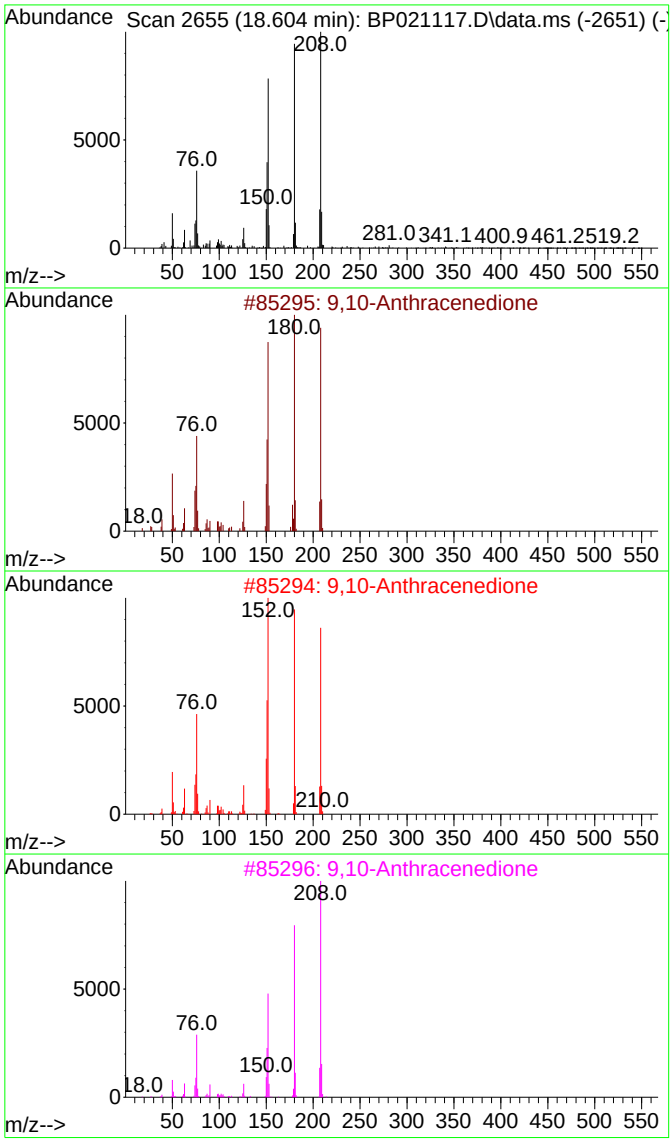
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

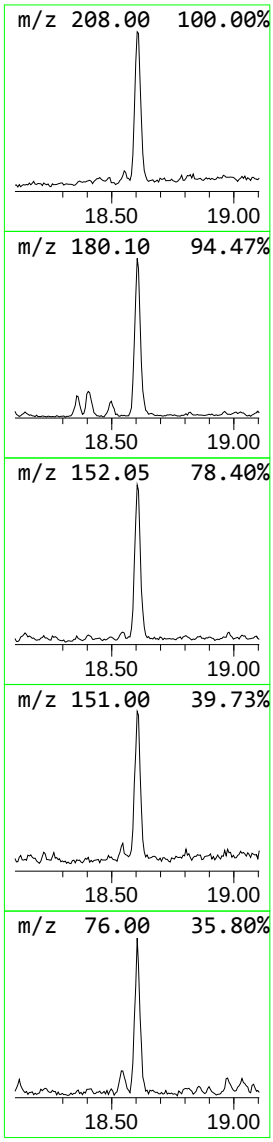
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 9,10-Anthracenedione Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.604	2.77 ng/ul	348939	Phenanthrene-d10	17.128

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
2	9,10-Anthracenedione	208	C14H8O2	000084-65-1	98
3	9,10-Anthracenedione	208	C14H8O2	000084-65-1	94
4	9,10-Anthracenedione	208	C14H8O2	000084-65-1	87
5	Benzo[c]cinnoline	180	C12H8N2	000230-17-1	58





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

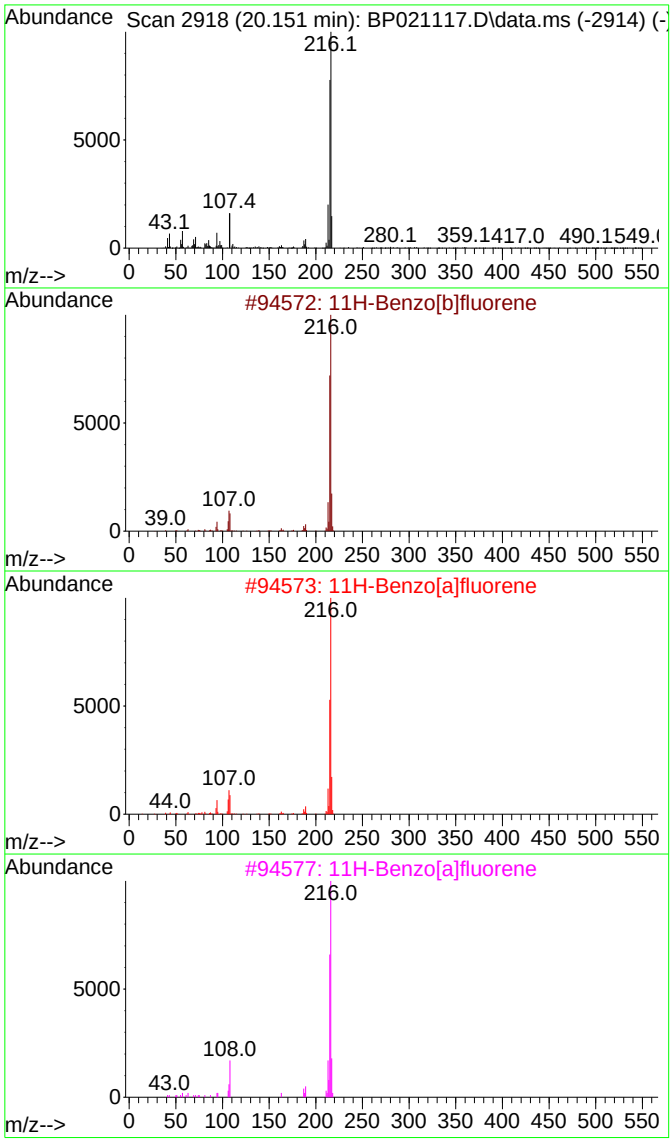
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

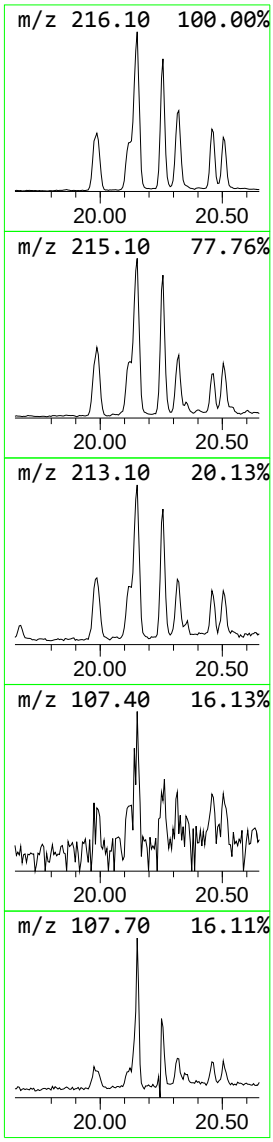
****

Peak Number 7 11H-Benzo[b]fluorene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.151	2.40 ng/ul	705378	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	93
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	80
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	80





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

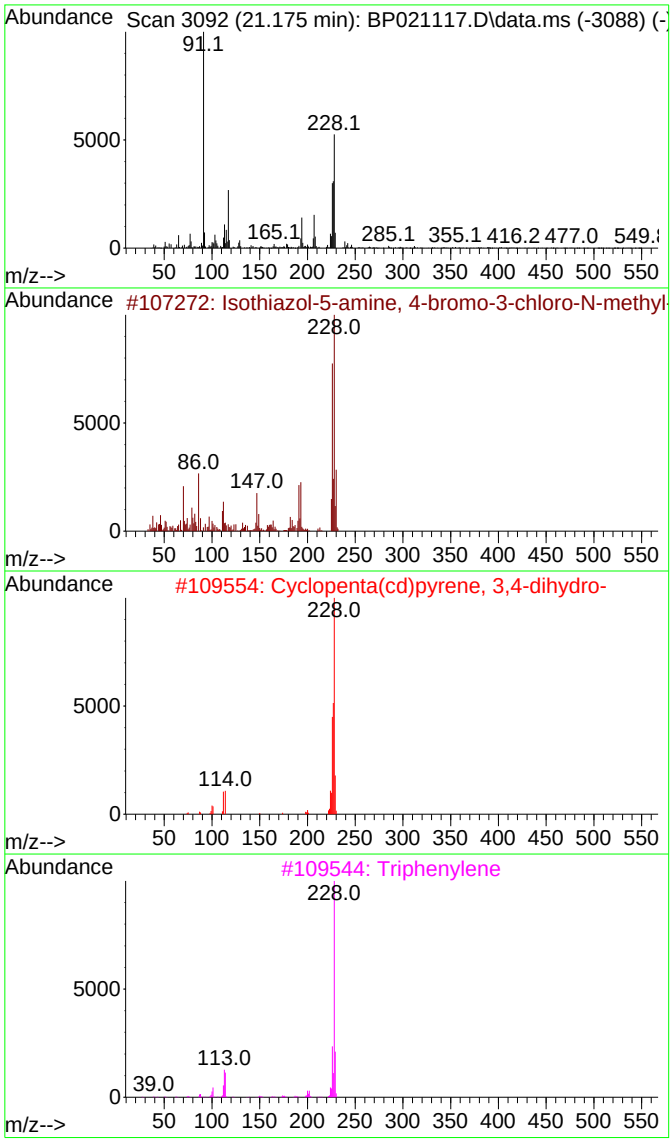
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

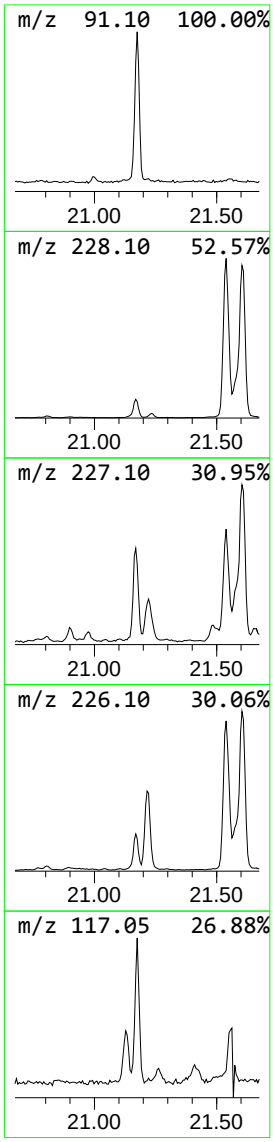
****

Peak Number 8 unknown-02 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.175	2.02 ng/ul	592126	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isothiazol-5-amine, 4-bromo-3-ch...	226	C4H4BrClN2S	1000272-59-9	35
2	Cyclopenta(cd)pyrene, 3,4-dihydro-	228	C18H12	025732-74-5	30
3	Triphenylene	228	C18H12	000217-59-4	30
4	Benzo[c]phenanthrene	228	C18H12	000195-19-7	27
5	Benzo[c]phenanthrene	228	C18H12	000195-19-7	25





Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

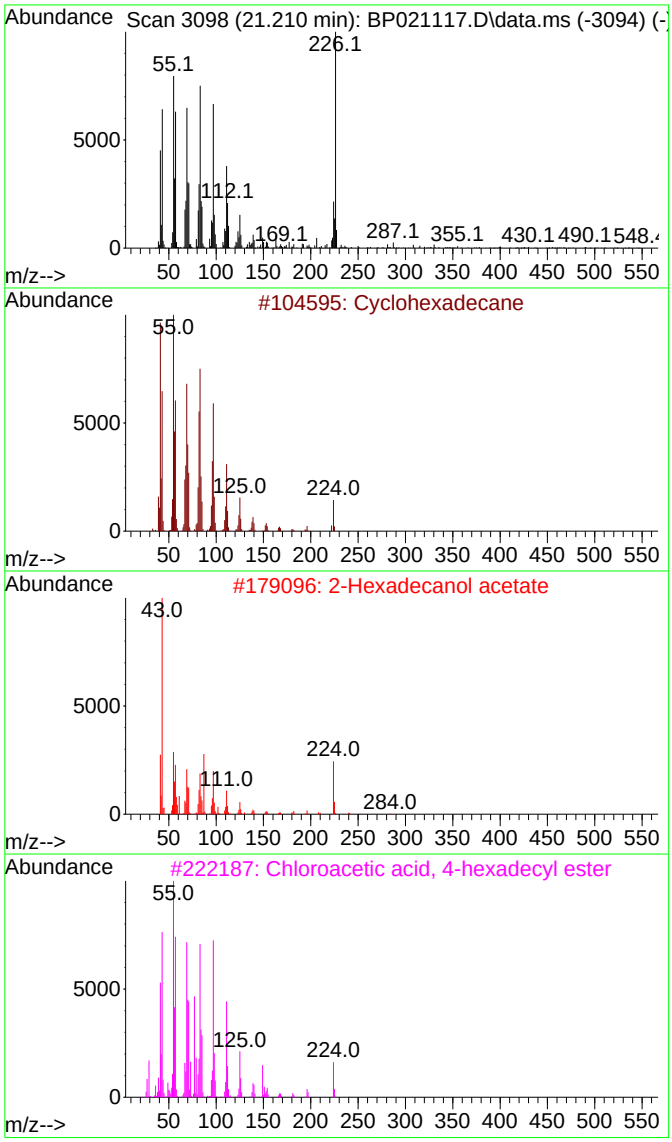
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

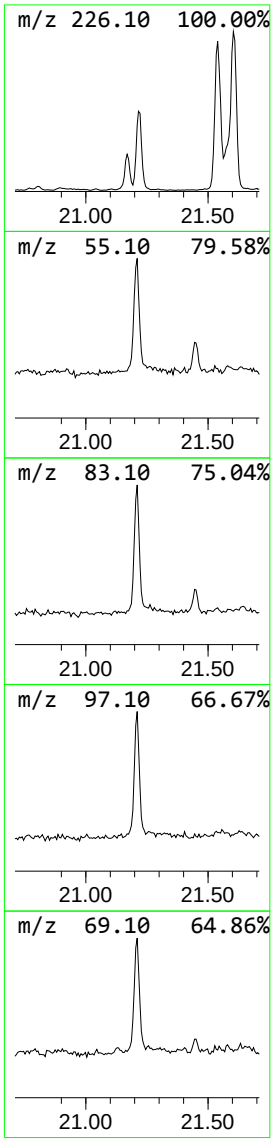
TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Cyclic21.210 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.210	4.59 ng/ul	1349820	Chrysene-d12	21.557

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexadecane	224	C16H32	000295-65-8	78
2	2-Hexadecanol acetate	284	C18H36O2	037590-88-8	45
3	Chloroacetic acid, 4-hexadecyl e...	318	C18H35ClO2	1000282-93-2	43
4	1-Hexadecanol	242	C16H34O	036653-82-4	38
5	Pentafluoropropionic acid, 4-hex...	388	C19H33F5O2	1000283-04-1	38





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072324\
Data File : BP021117.D
Acq On : 24 Jul 2024 04:00
Operator : MA/JU
Sample : P3238-16
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4BZ6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071024.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
1-Phenoxypropan...	11.187	4.0	ng/ul	296178	2	10.434	1462240	20.0
unknown-01	14.504	6.4	ng/ul	657352	3	14.310	2053570	20.0
n-Hexadecanoic ...	18.057	4.8	ng/ul	611012	4	17.128	2523820	20.0
Phenanthrene, 1...	18.087	3.0	ng/ul	372040	4	17.128	2523820	20.0
4H-Cyclopenta[d...	18.228	4.8	ng/ul	607449	4	17.128	2523820	20.0
9,10-Anthracene...	18.604	2.8	ng/ul	348939	4	17.128	2523820	20.0
11H-Benzo[b]flu...	20.151	2.4	ng/ul	705378	5	21.557	5876890	20.0
unknown-02	21.175	2.0	ng/ul	592126	5	21.557	5876890	20.0
(DEL) Alkane: C...	21.210	4.6	ng/ul	1349820	5	21.557	5876890	20.0